

# Classical Monte Carlo Simulations



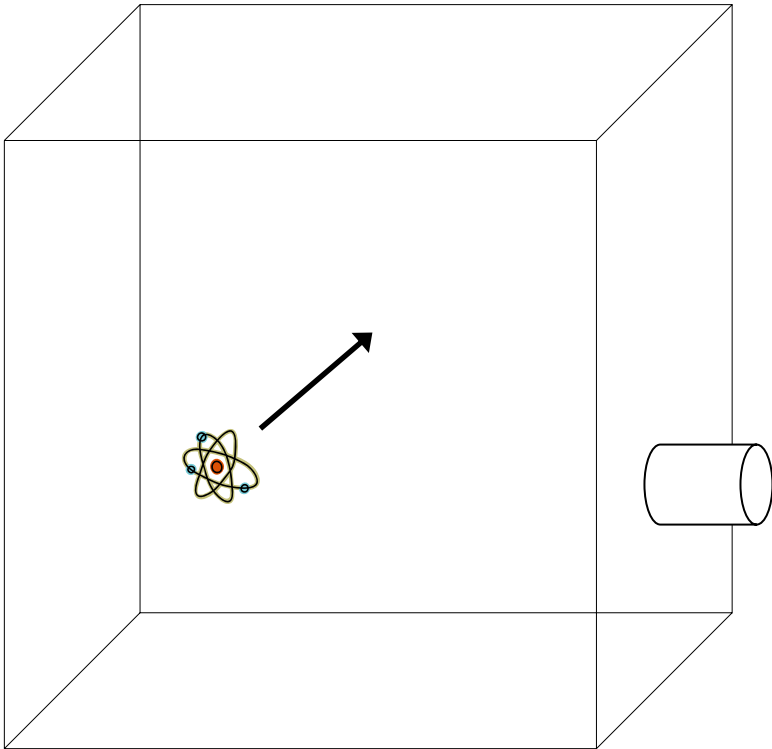
... you can calculate anything with dice.

# Introduction

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- the number of **bounces** (mean and variance)
- and **time** (mean and variance)

for a gas molecule at temperature  $T$  to leave this box:

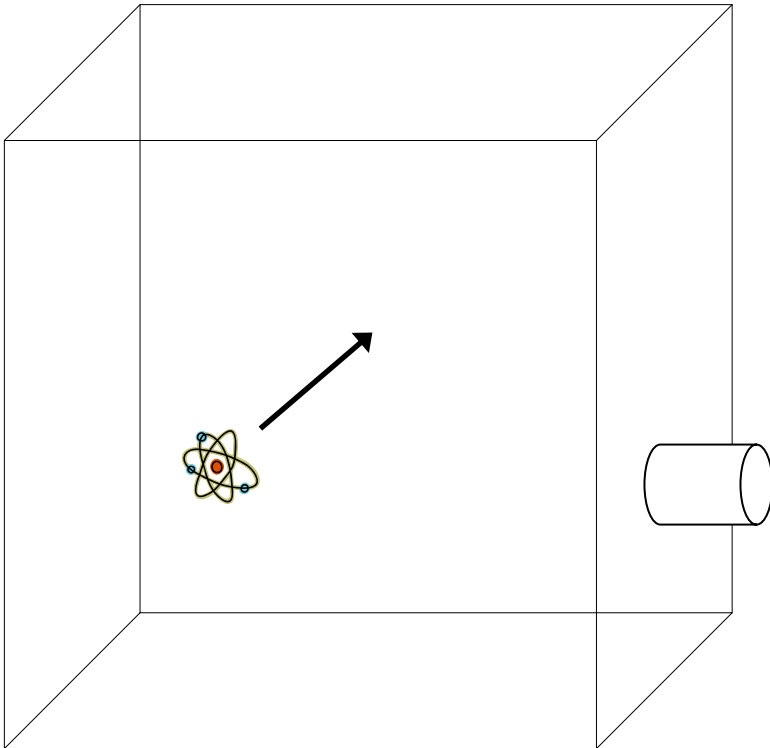


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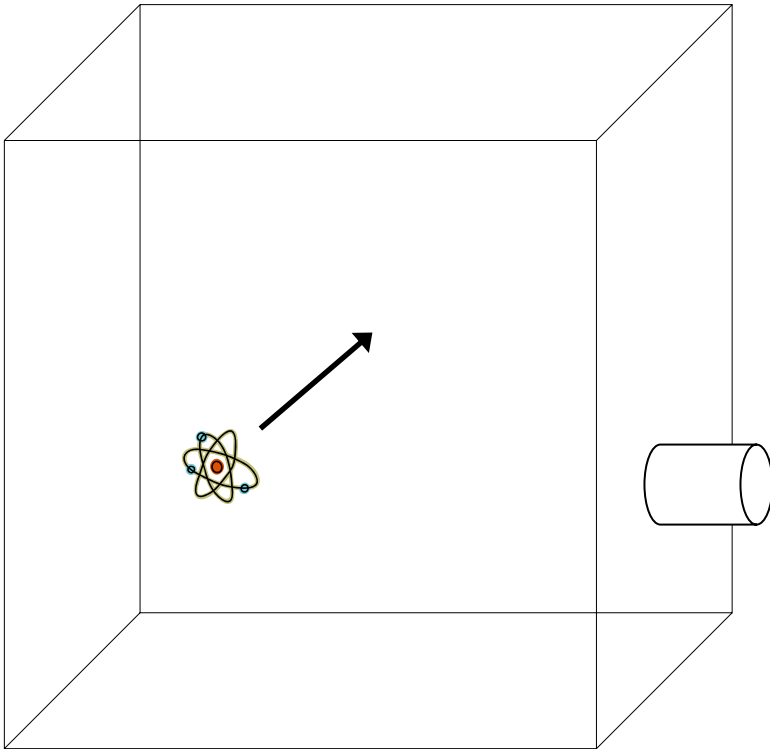
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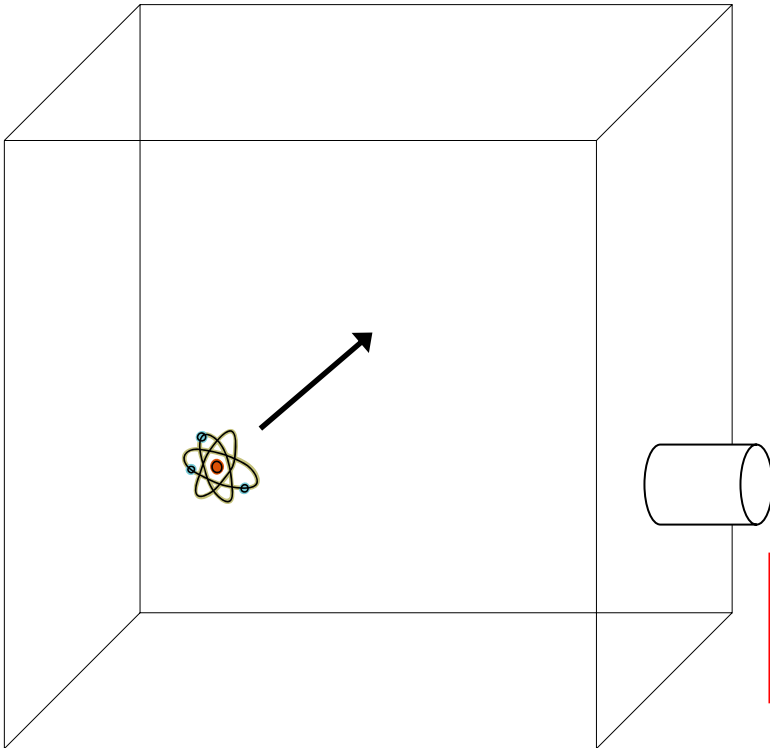
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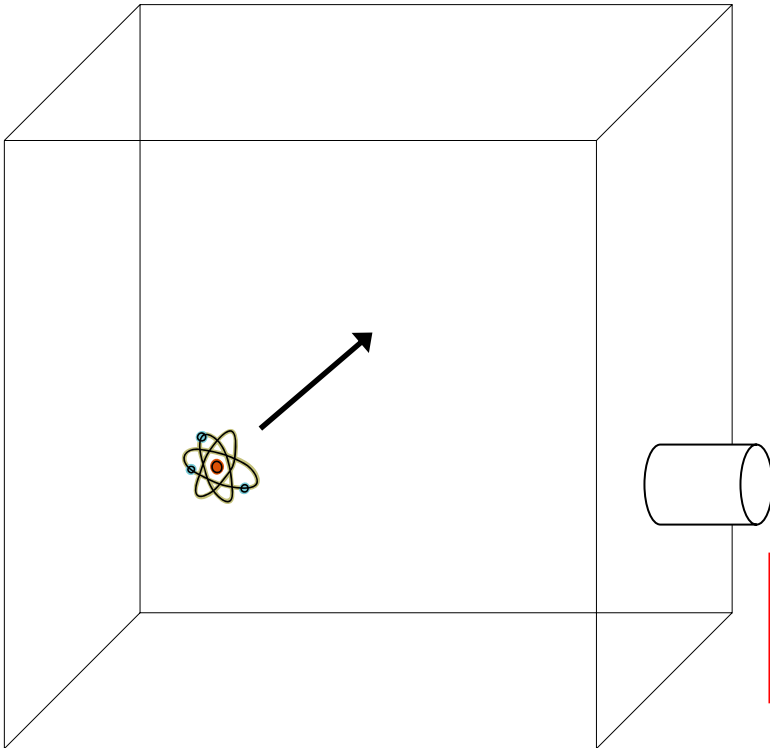
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... fairly simple and quick on a computer.

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# Definition

The Monte Carlo method is any numerical method in which the solution is obtained by ***averaging over many probabilistic simulation instances.***

# Example: Numerical Integration

The Monte Carlo method is frequently used to evaluate difficult integrals (in many dimensions):

$$\text{“calculus” average: } \langle f(x) \rangle_{[a,b]; \textit{calculus}} = \frac{1}{b-a} \int_a^b f(x) dx$$



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$x_i$ =probabilistic variable

i.e. choose  $x_i$ 's randomly on  $[a,b]$  with a uniform probability distribution.

# Theorem

If  $f(x)$  is well behaved on  $[a,b]$  (i.e. does not diverge), then in the limit of  $N \rightarrow \infty$ , the following is true (in the probabilistic sense)

$$\langle f(x) \rangle_{[a,b]; \text{calculus}} = \langle f(x) \rangle_{[a,b]; \text{statistical}, N} \pm \frac{\sigma_{f(x), N}}{\sqrt{N}}$$

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where 
$$\sigma_{f(x), N}^2 = \frac{1}{N-1} \left( \sum_{i=1}^N f(x_i)^2 - N \langle f(x) \rangle_{[a,b], N}^2 \right) = \frac{1}{N-1} \sum_{i=1}^N (f(x_i) - \langle f(x) \rangle)^2$$
  
= standard deviation of simulations

# Advantages

- Monte Carlo simulations are generally easy to formulate and set-up.
- Monte Carlo simulations are generally faster than other numerical methods, especially for problems in a large number of dimensions.